## **Amendments to the Specification:**

Please replace Table 1, on pages 19-21 with the following amended Table 1:

TABLE 1

SEQ NO:	ID Sequence (3' → 5')	Length (#bases)
21 22	CTAGAAGGAGAGATGGGTGCGAGAG AGAAGGAGAGAGAGGGUGCGAGAGCGUCAGUAUUAAGC	Target <sup>b</sup> Target <sup>b</sup>
1	CCCACGCTC	9
2	TTCCTCTCTA	12
3	CTTCCTCTCT	12
4	TCTTCCTCTC	12
5	TTCCTCTCTCTACCCACGCTC	21
6	CTTCCTCTCTGCCCACGCTC	22
7	TCTTCCTCTC <u>CG</u> CCCACGCTC	23
8	CTTCCTCTCTA	13
9	TTCCTCTCTA G G C	15 <del>15</del>
10	CTTCCTCTCT  G  G  C	15
11	CTTCCTCTCT  G G C C	16
12	CTTCCTCTCT  G  G  C  C  C	17
13	CTTCCTCTCT  G  G  C  C  C  G  G	19

TABLE	1	(continued)
TWDTP	1	(continued

SEQ ID	Deputition I addr.	
NO:	Sequence $(3' \rightarrow 5')$	Length (# bases)
14	CCCACGCTC C C G	12
15	CCCACGCTC C C G G	13
16	CCCACGCTC C C G G G	14
17	CCCACGCTC C C G G C C	16
18	CCCAC <u>T</u> CTC	9
19	CC <u>A</u> AC <u>T</u> CTC	9
20	TCTTCCTCTCTACCCACGCTCTC	25
23	$\mathtt{TTCCTCTCTACCCAC}_{\mathtt{T}}\mathtt{CTC}$	21
24	$\mathtt{TTCCTCTCTACC}$ $\mathtt{AAC}$ $\mathtt{TCTC}$	21
25	adamantane-CCCACGCTC	9
26	TTCCTCTCTA-cyclodextrin	12
27	CTTCCTCTCT-cyclodextrin	12
28	ATCTTCCTCT-cyclodextrin	<u>12</u>
29	CCCACGCTC C C adamantane-G	<u>12</u> <del>15</del>
30	CTCTTCCTCTCT G G C-cyclodextrin	<u>17</u>

<sup>\*</sup>underlined bases represent mismatches

<sup>&</sup>lt;sup>b</sup> sequence is  $5' \rightarrow 3'$ 

Please replace Table 2 on pages 24-25 with the following amended Table 2:

TABLE 2

Oligos (SEO ID NO:)	Complex <sup>a,b</sup>	Tm,°C
21 <del>1</del> 1	CTAGAAGGAGAGATGGGTGCGAGAG CCCACGCTC	49.1
21 <del>2</del> 2	CTAGAAGGAGAGATGGGTGCGAGAG TTCCTCTCTA	43.4
<u>21</u> 3 <u>3</u>	CTAGAAGGAGAGATGGGTGCGAGAG CTTCCTCTCT	43.6
<u>21</u> 4 <u>4</u>	CTAGAAGGAGAGATGGGTGCGAGAG TCTTCCTCTCTC	45.0
<u>21</u> <del>5</del> <u>5</u>	CTAGAAGGAGAGATGGGTGCGAGAG TTCCTCTCTACCCACGCTC	67.7
<u>21</u> <del>6</del> <u>6</u>	CTAGAAGGAGAGATGGGTGCGAGAG CTTCCTCTCT <u>G</u> CCCACGCTC	64.2
<u>21</u> 7 7	CTAGAAGGAGAGATGGGTGCGAGAG TCTTCCTCTC <u>CG</u> CCCACGCTC	59.9
21 <del>1+2</del> 1+2	CTAGAAGGAGAGATGGGTGCGAGAG TTCCTCTCTACCCACGCTC	47.8
<u>21</u> <del>1+3</del> 1+3	CTAGAAGGAGAGATGGGTGCGAGAG CTTCCTCTCT CCCACGCTC	44.4
21 <del>1+4</del> 1+4	CTAGAAGGAGAGATGGGTGCGAGAG TCTTCCTCTC CCCACGCTC	45.9
21 <sup>1+8</sup> 1+8	CTAGAAGGAGAGATGGGTGCGAGAG CTTCCTCTCTACCCACGCTC	50.5

a = underlined bases represent mismatches

Please replace Table 3 on pages 26-27 with the following amended Table 3:

TABLE 3

Oligos (SEQ ID NOS:)	Complex	Tm,°C
1100.7	COMPTCX	11117 C
21	CTAGAAGGAGAGATGGGTGCGAGAG	
10+14	CTTCCTCTCTCT CCCACGCTC	45.9
	G C	
	G C	
	C G	

b = The target sequence is bolded and is 5'to 3'; the oligo corresponding to SEO ID No. 1 is italicized

## TABLE 3 (continued)

Oligos		
(SEQ ID NOS:)	Complex <sup>a</sup>	Tm,°C
<u>21</u>	CTAGAAGGAGAGATGGGTGCGAGAG	
11+15	CTTCCTCTCT CCCACGCTC	47.3
	G C	
	G C	
	C G	
	C G	
<u>21</u>	CTAGAAGGAGAGATGGGTGCGAGAG	
12+16	CTTCCTCTCT CCCACGCTC	48.4
	G C	
	G C	
	C G	
	C G	
	G C	
<u>21</u>	CTAGAAGGAGAGATGGGTGCGAGAG	
13+1	CTTCCTCTCT CCCACGCTC	53.2
	G C	
	G C	
	C G	
	C G	
	G C C G	
	G C	
	G C	
<u>21</u>	CTAGAAGGAGAGATGGGTGCGAGAG	
9+14	TTCCTCTCTACCCACGCTC	47.9
	GC	
	GC	
	CG	

<sup>\*</sup> Target is bolded and is 5' to 3'; complementary cooperative oligonucleotides are 3' to 5'